

1. Combine finite difference approximations for $\partial u / \partial t$ at $x = x_i$

$$\left. \frac{\partial u}{\partial t} \right|_{t_k, x_i} = \frac{u_i^k - u_i^{k-1}}{\Delta t} + \mathcal{O}(\Delta t) \quad (1)$$

and $\partial^2 u / \partial x^2$ at time t_k

$$\left. \frac{\partial^2 u}{\partial x^2} \right|_{t_k, x_i} = \frac{u_{i-1}^k - 2u_i^k + u_{i+1}^k}{\Delta x^2} + \mathcal{O}(\Delta x^2). \quad (2)$$

to get

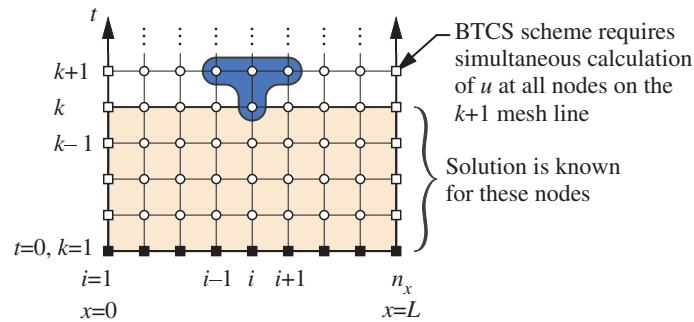
$$\frac{u_i^k - u_i^{k-1}}{\Delta t} = \alpha \frac{u_{i-1}^k - 2u_i^k + u_{i+1}^k}{\Delta x^2} + \mathcal{O}(\Delta t) + \mathcal{O}(\Delta x^2) \quad (3)$$

Drop the truncation error terms, shift the k subscripts by one, and move all u^{k+1} terms to the left hand side

$$\left[-\frac{\alpha}{\Delta x^2} \right] u_{i-1}^{k+1} + \left[\frac{1}{\Delta t} + \frac{2\alpha}{\Delta x^2} \right] u_i^{k+1} + \left[-\frac{\alpha}{\Delta x^2} \right] u_{i+1}^{k+1} = \frac{1}{\Delta t} u_i^k \quad (4)$$

Equation (4) is the computational formula for the BTCS scheme. It is an *implicit* scheme because all u^{k+1} values are coupled and must be updated simultaneously.

2. Computational Molecule



3. The BTCS method is *unconditionally stable* for the heat equation.

The benefit of stability comes at a cost of increased complexity of solving a linear system of equations at each time step.

4. For the one-dimensional heat equation, the linear system of equations for the BTCS method can be organized into a tridiagonal matrix.

$$\begin{bmatrix} a_1 & b_1 & 0 & 0 & 0 & 0 \\ c_2 & a_2 & b_2 & 0 & 0 & 0 \\ 0 & c_3 & a_3 & b_3 & 0 & 0 \\ 0 & 0 & \ddots & \ddots & \ddots & 0 \\ 0 & 0 & 0 & c_{n_x-1} & a_{n_x-1} & b_{n_x-1} \\ 0 & 0 & 0 & 0 & c_{n_x} & a_{n_x} \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \\ u_3 \\ \vdots \\ u_{n_x-1} \\ u_{n_x} \end{bmatrix} = \begin{bmatrix} d_1 \\ d_2 \\ d_3 \\ \vdots \\ d_{n_x-1} \\ d_{n_x} \end{bmatrix} \quad (5)$$

where the coefficients of the matrix are stored compactly as vectors

$$a_i = (1/\Delta t) + (2\alpha/\Delta x^2), \quad b_i = c_i = -\alpha/\Delta x^2, \quad d_i = (1/\Delta t)u_i^k. \quad (6)$$

This system of equations is efficiently solved with a form of LU factorization. The LU factors need to be computed only once before the first time step.

5. MATLAB implementation: code from `demoBTCS`

```
% --- Coefficients of the tridiagonal system
b = (-alfa/dx^2)*ones(nx,1); % Super diagonal: coefficients of u(i+1)
c = b; % Subdiagonal: coefficients of u(i-1)
a = (1/dt)*ones(nx,1) - (b+c); % Main Diagonal: coefficients of u(i)
a(1) = 1; b(1) = 0; % Fix coefficients of boundary nodes
a(end) = 1; c(end) = 0;
[e,f] = tridiagLU(a,b,c); % Save LU factorization

% --- Assign IC and save BC values in ub. IC creates u vector
x = linspace(0,L,nx)'; u = sin(pi*x/L); ub = [0 0];

% --- Loop over time steps
for k=2:nt
    d = [ub(1); u(2:nx-1)/dt; ub(2)]; % Update RHS, preserve BC
    u = tridiagLUSolve(e,f,b,d); % Solve the system
end
```

A more general implementation is in `heatBTCS`.

6. The `demoBTCS` function demonstrates the correct behavior of truncation error as Δx and Δt are reduced.

>> `convBTCS`

nx	nt	error	E(j)/E(j-1)	p
4	5	5.346e-02	0.0000	0.0000
8	21	1.186e-02	0.2219	2.1723
16	92	2.716e-03	0.2290	2.1268
32	386	6.522e-04	0.2401	2.0581
64	1589	1.594e-04	0.2444	2.0326
128	6453	3.939e-05	0.2471	2.0168
256	26012	9.790e-06	0.2485	2.0084
512	104452	2.440e-06	0.2493	2.0042

nx	nt	error	E(j)/E(j-1)	p
4	4	6.444e-02	0.0000	0.0000
8	8	2.737e-02	0.4248	1.2353
16	16	1.276e-02	0.4661	1.1014
32	32	6.172e-03	0.4838	1.0475
64	64	3.037e-03	0.4921	1.0230
128	128	1.507e-03	0.4961	1.0113
256	256	7.504e-04	0.4981	1.0056
512	512	3.745e-04	0.4990	1.0028

